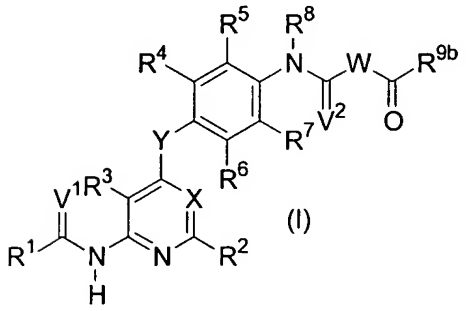
AMENDMENTS TO THE CLAIMS

1. (Original) A compound represented by the following formula, a salt thereof or a hydrate of the foregoing:



(I)

wherein R^1 represents C_{1-6} alkyl, C_{2-6} alkenyl, C_{2-6} alkynyl, C_{3-10} cycloalkyl, C_{6-10} aryl, C_{1-6} alkoxy, 5- to 10-membered heteroaryl, a 3- to 10-membered non-aromatic heterocyclic group or a group represented by the formula $-NR^{11a}R^{11b}$, and R^1 may be substituted with a substituent selected from Substituent Group A or Substituent Group B, wherein R^{11a} and R^{11b} may be the same or different and each represents hydrogen, C_{1-6} alkyl, C_{3-6} alkenyl, C_{3-6} alkynyl, C_{3-10} cycloalkyl, C_{6-10} aryl, C_{1-6} alkoxy, 5- to 10-membered heteroaryl or a 4- to 10-membered non-aromatic heterocyclic group, and R^{11a} and R^{11b} may be substituted with a substituent selected from Substituent Group A or Substituent Group B;

R^2 and R^3 represent hydrogen;

R^4, R^5, R^6 and R^7 may be the same or different and each represents hydrogen, halogen, hydroxyl, cyano, trifluoromethyl, C_{1-6} alkyl, C_{2-6} alkenyl, C_{2-6} alkynyl, C_{1-6} alkoxy, amino, mono-C_{1-6} alkylamino, di-C_{1-6} alkylamino or a group represented by the formula $-CO-R^{12}$, wherein R^{12} represents hydrogen, hydroxyl, C_{1-6} alkyl, C_{1-6} alkoxy, amino, mono-C_{1-6} alkylamino or di-C_{1-6} alkylamino;

R^8 represents hydrogen or C_{1-6} alkyl;

R^{9b} represents a 3- to 10-membered non-aromatic heterocyclic group (limited to a group having nitrogen as a ring constituent atom, the nitrogen having a bonding hand) or a group represented by the formula $-NR^{11a}R^{11b}$ wherein R^{11a} and R^{11b} represent the same meanings as recited above, and R^{9b} may be substituted with a substituent selected from Substituent Group A or Substituent Group B;

V^1 and V^2 may be the same or different and each represents oxygen or sulfur;

W represents a direct bond or a group represented by the formula $-C(R^{W1})(R^{W2})-$ wherein R^{W1} and R^{W2} are the same or different and each represents hydrogen, halogen, C_{1-6} alkyl or C_{1-6} alkoxy;

X represents a group represented by the formula $-C(R^{10})=$ or nitrogen, wherein R^{10} represents hydrogen, halogen, cyano, C_{1-6} alkyl, C_{2-6} alkenyl, C_{2-6} alkynyl or a group represented by the formula $-CO-R^{12}$, wherein R^{12} represents the same meaning as recited above; and

Y represents oxygen, sulfur, sulfinyl, sulfonyl or a group represented by the formula $-N(R^Y)-$, wherein R^Y represents hydrogen or C_{1-6} alkyl,

wherein Substituent Group A consists of halogen, hydroxyl, mercapto, nitro, cyano and oxo;

wherein Substituent Group B consists of C_{1-6} alkyl, C_{2-6} alkenyl, C_{2-6} alkynyl, C_{3-10} cycloalkyl, C_{6-10} aryl, 5- to 10-membered heteroaryl, a 3- to 10-membered non-aromatic heterocyclic group, C_{1-6} alkoxy, C_{3-6} alkenyloxy, C_{3-6} alkynyloxy, C_{3-10} cycloalkoxy, C_{6-10} aryloxy, 5- to 10-membered heteroaryloxy, 4- to 10-membered non-aromatic heterocyclicoxy, C_{1-6} alkylthio, C_{3-6} alkenylthio, C_{3-6} alkynylthio, C_{3-10} cycloalkylthio, C_{6-10} arylthio, 5- to 10-membered heteroarylthio, 4- to 10-membered non-aromatic heterocyclicthio and a group

represented by the formula $-T^1-T^2-T^3$, and each group in Substituent Group B may be substituted with a substituent selected from Substituent Group C, wherein T^1 represents a direct bond or C_{1-6} alkylene, T^2 represents carbonyl, sulfinyl, sulfonyl, a group represented by the formula -C(=O)-O-, a group represented by the formula -O-C(=O)-, a group represented by the formula $-SO_2-O-$, a group represented by the formula $-O-SO_2-$, a group represented by the formula $-NR^{T1}-$, a group represented by the formula $-C(=O)-NR^{T1}-$, a group represented by the formula $-NR^{T1}-C(=O)-$, a group represented by the formula $-SO_2-NR^{T1}-$ or a group represented by the formula $-NR^{T1}-SO_2-$, T^3 represents hydrogen, C_{1-6} alkyl, C_{3-6} alkenyl, C_{3-6} alkynyl, C_{3-10} cycloalkyl, C_{6-10} aryl, 5- to 10-membered heteroaryl or a 4- to 10-membered non-aromatic heterocyclic group, and R^{T1} represents hydrogen or C_{1-6} alkyl; and

wherein Substituent Group C consists of halogen, hydroxyl, mercapto, nitro, cyano, oxo, C_{1-6} alkyl, C_{2-6} alkenyl, C_{2-6} alkynyl, C_{3-10} cycloalkyl, C_{6-10} aryl, 5- to 10-membered heteroaryl, a 3- to 10-membered non-aromatic heterocyclic group, C_{1-6} alkoxy and C_{1-6} alkylthio.

2. (Original) A compound according to Claim 1, a salt thereof or a hydrate of the foregoing, wherein R^1 represents C_{1-6} alkyl optionally substituted with a substituent selected from Substituent Group A or Substituent Group B recited in Claim 1.

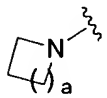
3. (Original) A compound according to Claim 1, a salt thereof or a hydrate of the foregoing, wherein R^1 represents C_{1-6} alkyl optionally substituted with a substituent selected from Substituent Group D,

wherein Substituent Group D consists of amino, mono-C_{1-6} alkylamino and di-C_{1-6} alkylamino.

4. (Original) A compound according to Claim 1, a salt thereof or a hydrate of the

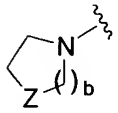
foregoing, wherein R^1 represents a 3- to 10-membered non-aromatic heterocyclic group optionally substituted with a substituent selected from Substituent Group A or Substituent Group B recited in Claim 1.

5. (Original) A compound according to Claim 1, a salt thereof or a hydrate of the foregoing, wherein R^1 represent a group represented by the formula (II):


(II)

wherein a represents an integer of 1 to 4,

or a group represented by the formula (III):


(III)

wherein b represents an integer of 1 to 3, and Z represents oxygen, sulfur, carbonyl, sulfonyl or a group represented by the formula $-NR^Z-$, wherein R^Z represents hydrogen or C_{1-6} alkyl,

and the groups represented by the formula (II) or (III) may be substituted with a substituent selected from Substituent Group A or Substituent Group B recited in Claim 1.

6. (Original) A compound according to Claim 1, a salt thereof or a hydrate of the foregoing, wherein R^1 represents azetidin-1-yl optionally substituted with a substituent selected from Substituent Group E, pyrrolidin-1-yl optionally substituted with a substituent selected from Substituent Group E, piperidin-1-yl optionally substituted with a substituent selected from Substituent Group E, azepan-1-yl optionally substituted with a substituent selected from Substituent Group E, piperazin-1-yl optionally substituted with a substituent selected from Substituent Group E, diazepan-1-yl optionally substituted with a substituent selected from

Substituent Group E, morpholin-4-yl optionally substituted with a substituent selected from Substituent Group E, thiomorpholin-4-yl optionally substituted with a substituent selected from Substituent Group E or 1,1-dioxothiomorpholin-4-yl optionally substituted with a substituent selected from Substituent Group E,

wherein Substituent Group E consists of halogen, hydroxyl, mercapto, cyano, formyl, oxo, C_{1-6} alkyl, C_{3-10} cycloalkyl, C_{1-6} alkoxy, amino, mono-C_{1-6} alkylamino, di-C_{1-6} alkylamino, azetidinyl, pyrrolidinyl, piperidinyl, piperazinyl, diazepanyl and a group represented by $-T^4-T^5$, wherein T^4 represents carbonyl or sulfonyl, and T^5 represents C_{1-6} alkyl, C_{3-10} cycloalkyl, azetidinyl, pyrrolidinyl, piperidinyl, hydroxyl, C_{1-6} alkoxy, amino, mono-C_{1-6} alkylamino or di-C_{1-6} alkylamino,

where each group included in Substituent Group E may be substituted with hydroxyl, C_{1-6} alkyl, di-C_{1-6} alkylamino, azetidinyl or pyrrolidinyl.

7. (Original) A compound according to Claim 1, a salt thereof or a hydrate of the foregoing, wherein R^1 represents azetidin-1-yl optionally substituted with a substituent selected from Substituent Group E', pyrrolidin-1-yl optionally substituted with a substituent selected from Substituent Group E', piperidin-1-yl optionally substituted with a substituent selected from Substituent Group E', piperazin-1-yl optionally substituted with a substituent selected from Substituent Group E', diazepan-1-yl optionally substituted with a substituent selected from Substituent Group E' or morpholin-4-yl optionally substituted with a substituent selected from Substituent Group E',

wherein Substituent Group E' consists of methyl, ethyl, dimethylamino, azetidinyl, pyrrolidinyl, piperidinyl and piperazinyl,

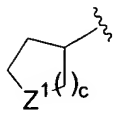
where each group included in Substituent Group E' may be substituted with hydroxyl, methyl, dimethylamino, azetidinyl or pyrrolidinyl.

8. (Original) A compound according to Claim 1, a salt thereof or a hydrate of the foregoing, wherein R^1 represents azetidin-1-yl optionally substituted with a substituent selected from Substituent Group E'", pyrrolidin-1-yl substituted with a substituent selected from Substituent Group E'" or piperidin-1-yl substituted with a substituent selected from Substituent Group E'"

wherein Substituent Group E'" consists of dimethylamino, azetidinyl, pyrrolidinyl, piperidinyl, dimethylaminomethyl, azetidin-1-ylmethyl, pyrrolidin-1-ylmethyl and piperidin-1-ylmethyl.

9. (Original) A compound according to Claim 1, a salt thereof or a hydrate of the foregoing, wherein R^1 represents a group represented by the formula $-NR^{11a}R^{11b}$, wherein R^{11a} and R^{11b} represent the same meaning as recited in Claim 1.

10. (Original) A compound according to Claim 1, a salt thereof or a hydrate of the foregoing, wherein R^1 represents a group represented by the formula $-NR^{11c}R^{11d}$, wherein R^{11c} represents hydrogen or C_{1-6} alkyl, and R^{11d} represents C_{1-6} alkyl or a group represented by the formula (IV):



(IV)

wherein c represents an integer of 1 to 3, and Z^1 represents oxygen, sulfur, carbonyl, sulfonyl or a group represented by the formula $-NR^{Z1}-$, wherein R^{Z1} represents hydrogen or C_{1-6} alkyl, and R^{11d} may be substituted with a substituent selected from Substituent Group A or Substituent

Group B recited in Claim 1.

11. (Original) A compound according to Claim 1, a salt thereof or a hydrate of the foregoing, wherein R^1 represents a group represented by the formula $-NR^{11e}R^{11f}$, wherein R^{11e} represents hydrogen or C_{1-6} alkyl, and R^{11f} represents C_{1-6} alkyl, pyrrolidin-3-yl, piperidin-3-yl, piperidin-4-yl or tetrahydropyran-4-yl, and R^{11f} may be substituted with a substituent selected from Substituent Group E recited in Claim 6.

12. (Original) A compound according to Claim 1, a salt thereof or a hydrate of the foregoing, wherein R^1 represents a group represented by the formula $-NR^{11g}R^{11h}$, wherein R^{11g} represents hydrogen or methyl, and R^{11h} represents n-propyl, n-butyl, pyrrolidin-3-yl, piperidin-3-yl, piperidin-4-yl or tetrahydropyran-4-yl, and R^{11h} may be substituted with a substituent selected from Substituent Group E",

wherein Substituent Group E" consists of methyl, ethyl, n-propyl, acetyl, dimethylamino, diethylamino, azetidinyl, pyrrolidinyl and piperazinyl,

where each group included in Substituent Group E" may be substituted with methyl or dimethylamino.

13. (Original) A compound according to Claim 1, a salt thereof or a hydrate of the foregoing, wherein R^1 represents a group represented by the formula $-N(CH_3)R^{11i}$, wherein R^{11i} represents n-propyl, n-butyl, pyrrolidin-3-yl or piperidin-4-yl, and R^{11i} is substituted with a substituent selected from Substituent Group E"",

wherein Substituent Group E"" consists of dimethylamino, diethylamino, dimethylaminoethyl, dimethylaminopropyl and 1-methylazetidin-3-yl.

14. (Currently amended) A compound according to ~~any one of Claims 1 to 13~~ Claim 1, a

salt thereof or a hydrate of the foregoing, wherein R^4, R^5, R^6 and R^7 may be the same or different and each represents hydrogen, halogen or C_{1-6} alkyl.

15 (Currently amended) A compound according to ~~any one of Claims 1 to 14~~ Claim 1, a salt thereof or a hydrate of the foregoing, wherein R^8 represents hydrogen.

16. (Currently amended) A compound according to ~~any one of Claims 1 to 15~~ Claim 1, a salt thereof or a hydrate of the foregoing, wherein V^1 represents oxygen.

17. (Currently amended) A compound according to ~~any one of Claims 1 to 16~~ Claim 1, a salt thereof or a hydrate of the foregoing, wherein X represents a group represented by the formula $-C(R^{10a})=$, wherein R^{10a} represents hydrogen, halogen or cyano.

18. (Currently amended) A compound according to ~~any one of Claims 1 to 16~~ Claim 1, a salt thereof or a hydrate of the foregoing, wherein X represents nitrogen.

19 (Currently amended) A compound according to ~~any one of Claims 1 to 18~~ Claim 1, a salt thereof or a hydrate of the foregoing, wherein Y represents oxygen.

20. (Currently amended) A compound according to ~~any one of Claims 1 to 19~~ Claim 1, a salt thereof or a hydrate of the foregoing, wherein W represents a group represented by the formula $-C(R^{W1})(R^{W2})-$ wherein R^{W1} and R^{W2} represent the same meanings as recited in Claim 1 and V^2 represents oxygen.

21. (Currently amended) A compound according to ~~any one of Claims 1 to 19~~ Claim 1, a salt thereof or a hydrate of the foregoing, wherein W represents a group represented by the formula $-CH_2-$ and V^2 represents oxygen.

22. (Currently amended) A compound according to ~~any one of Claims 1 to 21~~ Claim 1, a salt thereof or a hydrate of the foregoing, wherein R^{9b} represents mono-C_{1-6} alkylamino

optionally substituted with a substituent selected from Substituent Group A or Substituent Group B recited in Claim 1, mono-C_{3-10} cycloalkylamino optionally substituted with a substituent selected from Substituent Group A or Substituent Group B recited in Claim 1, mono-C_{6-10} arylamino optionally substituted with a substituent selected from Substituent Group A or Substituent Group B recited in Claim 1, mono-5- to 10-membered heteroarylamino optionally substituted with a substituent selected from Substituent Group A or Substituent Group B recited in Claim 1 or mono-4- to 10-membered non-aromatic heterocyclic amino optionally substituted with a substituent selected from Substituent Group A or Substituent Group B recited in Claim 1.

23. (Currently amended) A compound according to ~~any one of Claims 1 to 21~~ Claim 1, a salt thereof or a hydrate of the foregoing, wherein R^{9b} represents mono-C_{3-10} cycloalkylamino optionally substituted with a substituent selected from Substituent Group A or Substituent Group B recited in Claim 1 or mono-C_{6-10} arylamino optionally substituted with a substituent selected from Substituent Group A or Substituent Group B recited in Claim 1.

24. (Original) A pharmaceutical composition comprising a compound according to claim 1, a salt thereof or a hydrate of the foregoing.

25. (Original) An inhibitor for hepatocyte growth factor receptor, comprising a compound according to Claim 1, a salt thereof or a hydrate of the foregoing.

26. (Original) An angiogenesis inhibitor comprising a compound according to Claim 1, a salt thereof or a hydrate of the foregoing.

27. (Original) An anti-tumor agent comprising a compound according to Claim 1, a salt thereof or a hydrate of the foregoing.

28. (Original) An anti-tumor agent according to Claim 27, wherein tumor is a pancreatic

cancer, a gastric cancer, a colorectal cancer, a breast cancer, a prostate cancer, a lung cancer, a renal cancer, a brain tumor or an ovarian cancer.

29. (Original) An inhibitor for cancer metastasis, comprising a compound according to Claim 1, a salt thereof or a hydrate of the foregoing.